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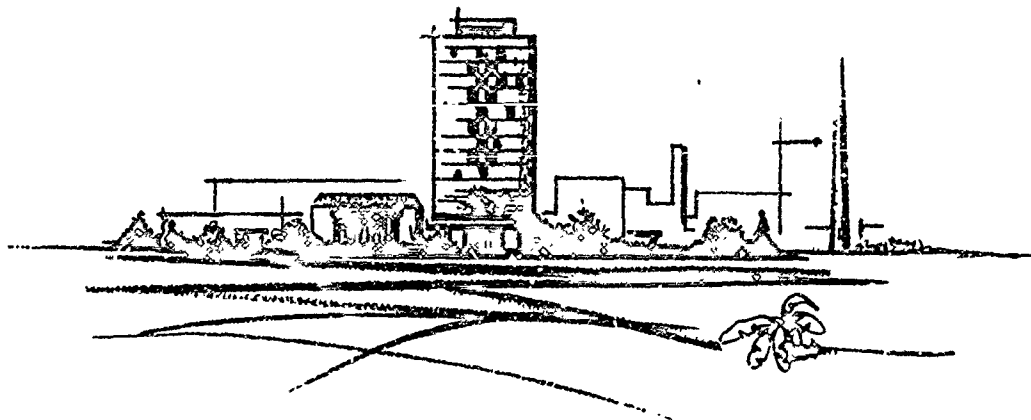
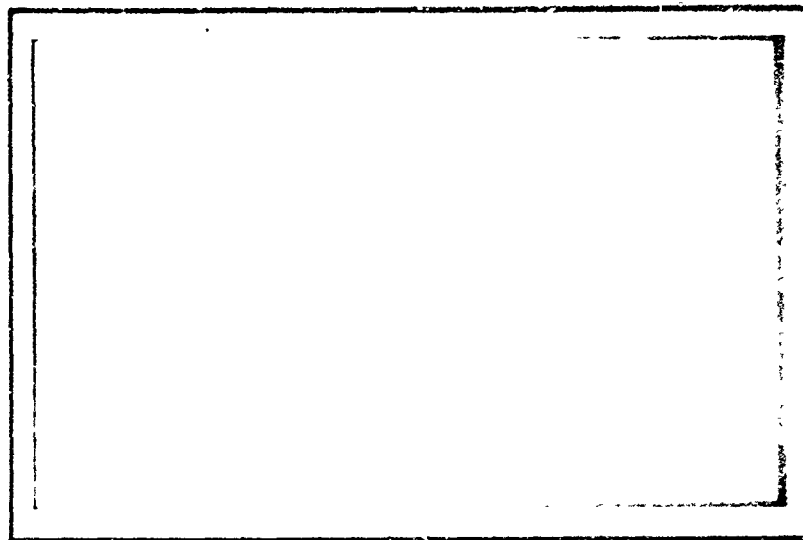
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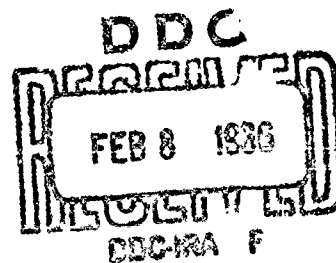
# RESEARCH REPORT



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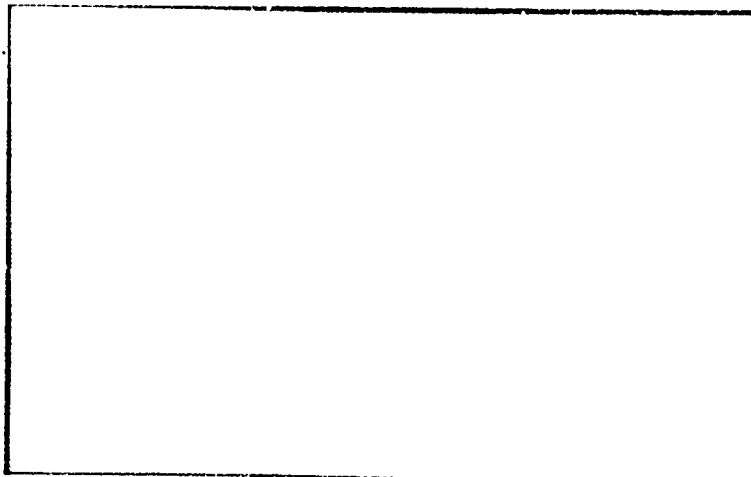
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⑨ FINAL REPORT.

⑥ HIGH-TEMPERATURE PROPERTIES AND  
ALLOYING BEHAVIOR OF THE REFRACTORY  
PLATINUM-GROUP METALS.

⑮ Contract (Nonr-2547(00)), ⑫ NR-039-067

to

OFFICE OF NAVAL RESEARCH

December 17, 1965

⑪ 17 Dec 65,

⑫ 4 p.

⑩ <sup>ter</sup>  
P. S. Rudman.  
Metal Science Group

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December 17, 1965

Dr. W. G. Rauch  
Acting Head, Metallurgy Branch  
Office of Naval Research  
Department of the Navy  
Washington, D.C. 20025

Dear Dr. Rauch:

Contract Nonr-2547(00), NR 039-067

Enclosed are two copies of the Final Report on the project, "High-Temperature Properties and Alloying Behavior of the Refractory Platinum-Group Metals".

Please let me know if you have any questions or comments concerning the information in this report.

Very truly yours,



Peter S. Rudman  
Fellow  
Metal Science Group

PSR:tam  
In duplicate  
Enc. (2)

cc: Mr. Edward P. Shute (2)  
ONR Resident Representative

# HIGH-TEMPERATURE PROPERTIES AND ALLOYING BEHAVIOR OF THE REFRACTORY PLATINUM-GROUP METALS

by

P. S. Rudman  
Metal Science Group

↓

The occurrence of many phases in platinum-group metal alloys, or even more generally in transition metal alloys, correlates amazingly simply with electron/atom ratio or as we have preferred to call it, group number. One example is the HCP structure that occurs in alloys in the average group number range 7-8.5. The axial ratio,  $c/a$ , ~~has been found to be~~ a sensitive measure of the electronic state. ~~We have determined the~~  $c/a$ -composition relationship in some 20 HCP alloys containing platinum-group metals. ~~The~~ *was determined,* axial ratio appears to correlate well with phase stability: the smaller the axial ratio, the more stable the phase, with the HCP phase becoming unstable relative to cubic phases as  $c/a$  ~~approaches~~ *approaches* 1.61.

However, group number is not the only structure determining factor. Atomic size is also very important. The Laves phases appeared to be examples of where atomic size plays an important structure determining role. A theoretical study of Laves phases based on an elastic model was performed. This study appears to furnish some insight into the origin of Laves structures and sets the groundwork for a study of their stability.

Before we can hope to understand the structural changes that occur on alloying, we surely must understand the origin of allotropism in pure metals. Accordingly, ~~a~~ *was* theoretical study of allotropism was initiated and is continuing. This study ~~has been~~ *was* very productive in providing clues to the phase-stabilizing factors. It has been tentatively concluded that the low temperature phases are generally characterized by a high density of states at the Fermi level and that the high temperature phases are characterized by high vibrational entropies.

↑

Publications in 1965

P. S. Rudman, "The Atomic Volumes of the Metallic Elements", Trans. AIME, 233, 864-872 (1965).

P. S. Rudman, "Atomic Volume in Laves Phases: A Hemisubstitutional Solid-Solution Elastic Model", Trans. AIME, 233, 872-878 (1965).

P. S. Rudman, "Lattice Parameters of Tantalum-Osmium Alloys", J. Less-Common Metals, 9, 77-79 (1965).